AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions and listings of claims in the application:

LISTING OF CLAIMS:

 (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof-or a hydrate thereof, represented by formula [I]

wherein,

$$\mathbb{R}^{d}$$

(wherein R^d is the same as described above) or a group represented by formula fiil; or,

in the case where either R^1 or R^2 represents a hydrogen atom, the other represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{4-10} alkyl group substituted by one or two aryl groups, a hydroxy C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkoxy C_{1-10} alkyl group, a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are the same as described above), a group represented by formula- $CHR^cOC(O)ZR^d$ (wherein Z, R^c and R^d are the same as described above), a group represented by formula $CHR^cOC(O)ZR^d$

$$\mathbb{R}^{d}$$

(wherein R^d is the same as described above) or a group represented by formula [ii];

X represents a hydrogen atom or a fluorine atom; and

Y represents -OCHR³R⁴, -SR³, -S(O)_nR⁵, -SCHR³R⁴, -S(O)_nCHR³R⁴, -NHCHR³R⁴,
-N(CHR³R⁴)(CHR³'R⁴), -NHCOR³ or -OCOR⁵ (wherein R³, R³', R⁴ and R⁴' are identical or different, and each represents a hydrogen atom, a C₁₋₁₀alkyl group, <u>or</u> a C₁₋₁₀alkenyl group, a phenyl group, a naphthyl group substituted by one to seven halogen atoms, a heteroaromatic group or a phenyl group substituted by one to five substituents selected from a group consisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxyearbonyl group, an amino group, a nitro group, a eyano group and a phenoxy group; R⁵ represents a C₁₋₁₀alkyl group, <u>or</u> a C₁₋₁₀alkenyl group, a phenyl group, a naphthyl group substituted by one to seven halogen atoms, a heteroaromatic group or a phenyl group substituted by one to five substituents selected from a group consisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a evano group and a phenoxy group; and n represents integer 1 or 2).

 (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof-or a hydrate thereof, represented by formula [II]

fwherein.

 R^1 and R^2 are identical or different, and each represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{2-10} alkyl group, a C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group), a group represented by formula- $C(C)CO)ZR^d$ (wherein $C(C)CO)ZR^d$) a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$) a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a C_{1-10} alkyl group, $C(C)CO)ZR^d$) and C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group, $C(C)CO)ZR^d$ (wherein a hydrogen atom, a C_{1-10} alkyl group,

$$\mathbb{R}^{d}$$

(wherein R^d is the same as described above) or a group represented by formula [ii]; or,

in the case where either R¹ or R² represents a hydrogen atom, the other represents a C₁.

10alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a-C₁₋₁₀alkyl group substituted by one or two-aryl-groups, a hydroxyC₂₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^{a-}and R^b are the same as described above), or a group represented by formula-CHR^cOC(O)ZR^d (wherein Z, R^c and R^d are the same as described above), a group represented by formula-CHR^cDC(O)ZR^d (wherein Z, R^c and R^d are the same as

(wherein R^d is the same as described above) or a group represented by formula [ii];

X represents a hydrogen atom or a fluorine atom; and

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Y represents -OCHR²R⁴, -SR³, -S(O)_nh⁵, -SCHR³R⁴, -S(O)_nCHR²R⁴, -NHCHR³R⁴,
-N(CHR³R⁴)(CHR³'R⁴), -NHCOR³ or -OCOR⁵ (wherein R³, R³', R⁴ and R⁴' are identical or different, and each represents a hydrogen atom, a C₁₋₁₀alkyl group, <u>or</u> a C₁₋₁₀alkenyl group, <u>a</u> naphthyl group, a naphthyl group substituted by one to seven halogen atoms, a heteroaromatic group or a phenyl group substituted by one to five substituents selected from a group consisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, <u>or</u> a C₁₋₁₀alkoxy group, a eyano group and a phenoxy group; R⁵ represents a C₁₋₁₀alkyl group, <u>or</u> a C₁₋₁₀alkenyl group, a henyl group, a naphthyl group, a naphthyl group, a naphthyl group, a naphthyl group substitute by one to seven halogen atoms, a heteroaromatic group or a phenyl group substituted by one to five substituents selected from a group consisting of a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a evano group and a phenoxy group; and n represents integer 1 or 24.

 (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof-or-a hydrate thereof according to claim 2, wherein in the formula [II],

R¹ and R² are identical or different, and each represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a C₁₋₁₀alkyl group substituted by one or two phenyl groups, a hydroxyC₂₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkyl group or a C₁₋₁₀alkyl group; or,

in the case where either R¹ or R² represents a hydrogen atom, the other represents a C₁.

10 alkyl group, a C₂₋₁₀ alkenyl group, a C₂₋₁₀ alkynyl group, a C₁₋₁₀ alkyl group substituted by one or

two phenyl groups, a hydroxyC₂₋₁₀ alkyl group, a halogenoC₁₋₁₀ alkyl group, an azidoC₁₋₁₀ alkyl

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group, an amino $C_{2\cdot 10}$ alkyl group, a $C_{1\cdot 10}$ alkoxy $C_{1\cdot 10}$ alkyl group or a $-C_{1\cdot 10}$ alkoxycarbonyl $C_{1\cdot 10}$ alkyl group.

 (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim
 wherein in the formula [II],

 R^1 and R^2 are identical or different, and each represents a C_{1-10} alkyl group, a C_{2-6} alkenyl group, a C_{2-6} alkynyl group, a C_{2-6} alkyl group substituted by one or two phenyl groups; a hydroxy C_{2-6} alkyl group, a halogeno C_{1-6} alkyl group, an azido C_{1-6} alkyl group, an azido C_{1-6} alkyl group, a C_{1-6} alkyl group or a C_{1-6} alkoxy C_{1-6} alkyl group or a C_{1-6} alkoxy C_{1-6} alkyl group or a C_{1-6} alkyl group; or,

in the case where either R¹ or R² represents a hydrogen atom, the other represents a C₁6alkyl group, a C₂₆alkenyl group, a C₂₆alkynyl group, a C₁₆alkyl group substituted by one or
two-phenyl groups, a hydroxyC₂₆alkyl group, a halogenoC₁₆alkyl group, an azidoC₁₆alkyl
group, an aminoC₂₆alkyl group, a C₁₆alkoxyC₁₆alkyl group or a C₁₆alkoxycarbonylC₁₆alkyl
group.

 (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II],

R¹ and R² are identical or different, and each represents a farnesyl group, a-C₁₋₁₀alkyl group substituted by one or two aryl groups, a C₁₋₁₀alkoxycarbonylC₁₋₁₀alkyl group, a 4morpholinylC₁₋₁₀alkyl group, a C₁₋₁₀alkyl group substituted by a group represented by formulaC(O)NR^aR^b (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group), a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a

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 C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group; and R^d represents a C_{1-10} alkyl group; or a C_{2-10} alkenyl group-or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$

(wherein R^d is the same as described above) or a group represented by formula [ii]; or,

in the case where either R^1 or R^2 represents a hydrogen atom, the other represents a farnesyl group, a C_{1-10} alkyl group substituted by one or two aryl groups, a C_{1-10} alkyl group, a C_{1-10} alkyl group, a C_{1-10} alkyl group, a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are the same as described above), or a group represented by formula- $C(O)ZR^d$ (wherein Z, R^c and R^d are the same as described above), a group represented by formula- $C(R^cO)ZR^d$ (wherein R^a) are the same as described above), a group represented by formula $C(R^d)$ formula $C(R^d$

$$\mathbb{R}^{d}$$

(wherein R^d is the same as described above) or a group represented by formula [ii]

6. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II],

R¹ and R² are identical or different, and each represents a farnesyl group, a-C₁₋₆alkyl group substituted by one or two aryl groups, a C₁₋₆alkoxycarbonylC₁₋₆alkyl group, a 4morpholinylC₁₋₆alkyl group, a C₁₋₆alkyl group substituted by a group represented by formulaC(O)NR³R³ (wherein R³ and R³ are identical or different, and each represents a hydrogen atom or a C₁₋₆alkyl group), a group represented by formula-CHR°OC(O)ZR⁴ (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R⁵ represents a hydrogen atom, a C₁₋₆alkyl group, or a C₂₋₆alkenyl group or an aryl group; and R⁴ represents a C₁₋₆alkyl group, or a C₂₋₆alkenyl group or an aryl group), a group represented by formula fil

$$\mathbb{R}^{d}$$
 \mathbb{O}

(wherein R^d is the same as described above) or a group represented by formula [ii]; or,

in the case where either R^1 or R^2 represents a hydrogen atom, the other represents a farnesyl group, a C_{L6} alkyl group substituted by one or two aryl groups, a C_{L6} alkoxycarbonyl C_{L6} alkyl group, a C_{L6} alkyl group, a C_{L6} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are the same as described above), or a group represented by formula- $CHR^cOC(O)ZR^d$ (wherein Z_1R^c and R^d are the same as described above), a group represented by formula fit

$$\mathbb{R}^{d}$$
 \mathbb{O}

(wherein R^d is the same as described above) or a group represented by formula [ii]

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7. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester

derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

2, wherein in the formula [II], R² represents a hydrogen atom.

8. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester

derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

2. wherein in the formula [III, R² represents a hydrogen atom; and X represents a fluorine atom.

9. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester

derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

2, wherein in the formula [II], wherein R² represents a hydrogen atom; and X represents a

hydrogen atom.

10. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester

derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

2, wherein in the formula [II], R^2 represents a hydrogen atom; X represents a fluorine atom; and

Y represents -OCHR³R⁴ (wherein R³ and R⁴ are the same as described above).

11. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester

derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim

2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and

Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above).

12. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester

derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim

2, wherein in the formula [II], \mathbb{R}^2 represents a hydrogen atom; X represents a fluorine atom; and

Y represents -SR3 (wherein R3 is the same as described above).

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13. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -S(O)₀CHR³R⁴(wherein R³, R⁴ and n are the same as described above).

- 14. (currently amended): A 2-amino-bicyclo[3.1.0]hexano-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- 15. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; and Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³', R⁴ and R⁴' are the same as described above).
- 16. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], represents a hydrogen atom; X represents a hydrogen atom; and Y represents -OCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- 17. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], represents a hydrogen atom; X represents a hydrogen atom; and Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim

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wherein in the formula [II], represents a hydrogen atom; X represents a hydrogen atom; and Y represents -SR³ (wherein R³ is the same as described above).

- 19. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; and Y represents -S(O)_nCHR³R⁴ (wherein R³, R⁴ and n are the same as described above).
- 20. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], wherein R² represents a hydrogen atom; X represents a hydrogen atom; and Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above).
- 21. (currently amended): A 2-amino-bicyclo[3.1.0]hexanc-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; and Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³', R⁴ and R⁴' are the same as described above).
- 22. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a-C₁₋₁₀alkyl group, a balogenoC₁₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkyl gro

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 $_{10}$ alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group).

23. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

R¹ represents a group represented by formula-CHR COC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C₁₋₁₀alkyl group, or a C₂₋₁₀alkenyl group-or an aryl group; and R^d represents a C₁₋₁₀alkyl group; or a C₂₋₁₀alkenyl group or an aryl group, a group represented by formula [i]

$$\mathbb{R}^{d}$$

(wherein R⁴ is the same as described above) or a group represented by formula [ii]

24. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative or a pharmaceutically acceptable salt thereof or a hydrate-thereof according to claim

2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, a nazido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

25. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR $^{c}OC(O)ZR^{d}$ (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^{c} represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group; and R^{d} represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group, a group represented by formula fill

$$\mathbb{R}^{d}$$
 0

(wherein R^{-th}is the same as described above) or a group represented by formula [ii]

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26. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents-SR³ (wherein R³ is the same as described above); and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, a nazidoC₁₋₁₀alkyl group, a naminoC₂₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxyCarbonylC₁₋₁₀alkyl group, a farnesyl group, a 4 morpholinylC₁₋₁₀alkyl group or a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group).

27. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -SR³ (wherein R³ is the same as described above); and

 R^1 represents a group represented by formula-CHR $^{\circ}$ OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 \mathbb{O}

(wherein R⁴ is the same as described above) or a group represented by formula [ii]

28. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -S(O)_nCHR³R⁴ (wherein R³, R⁴ and n are the same as described above); and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a C₁₋₁₀alkyl group substituted by one or two aryl groups, a hydroxyC₂₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxyCarbonylC₁₋₁₀alkyl group, a farnesyl group, a 4-morpholinylC₁₋₁₀alkyl group or a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group).

 (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim

2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -S(O)_nCHR³R⁴ (wherein R³,R⁴ and n are the same as described above); and

 R^1 represents a group represented by formula-CHR $^{\circ}$ OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, a C_{1-10} alkyl group, \underline{or} a C_{2-10} alkenyl group or an aryl group; and R^d represents a C_{1-10} alkyl group, \underline{or} a C_{2-10} alkenyl group or an aryl group, a group represented by formula [i]

$$\mathbb{R}^{d}$$
 \mathbb{O}

(wherein R^d is the same as described above) or a group represented by formula [ii]

30. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above): and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a C₁₋₁₀alkyl group substituted by one or two aryl groups, a hydroxyC₂₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, a C₁₋₁₀alkyl group, a

₁₀alkoxycarbonylC₁₋₁₀alkyl group, a farnesyl group, a 4 morpholinylC₁₋₁₀alkyl group or a C₁.

₁₀alkyl group substituted by a group represented by formula-C(O)NR*R^b (wherein R* and R^b are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group).

31. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR c OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, \underline{or} a C_{2-10} alkenyl group- \underline{or} and R^d represents a C_{1-10} alkyl group- \underline{or} a C_{2-10} alkenyl group- \underline{or} and R^d represents a C_{1-10} alkyl group- \underline{or} a C_{2-10} alkenyl group- \underline{or} and R^d represented by formula [i]

(wherein R^d is the same as described above) or a group represented by formula [ii]

(currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester
 derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim

2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³', R⁴ and R⁴' are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, a nazido C_{1-10} alkyl group, a namino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

33. (currently amended): A 2-amino-bicyclo[3.1.0]hexanc-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³, R⁴ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR $^{\circ}$ OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or-an-aryl group; and R^d represents a C_{1-10} alkyl group; or a C_{2-10} alkenyl group-or-an-aryl-group), a group represented by formula [i]

$$\mathbb{R}^{d}$$
 0 0

(wherein R^d is the same as described above) or a group represented by formula [ii]

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34. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents-OCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

35. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents-OCHR²R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a group represented by formula-CHR $^{\circ}$ OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, a C_{1-10} alkyl group, \underline{or} a C_{2-10} alkenyl group- \underline{or} an \underline{aryl} group; and R^d represents a C_{1-10} alkyl group, \underline{or} a C_{2-10} alkenyl group, \underline{or} a C_{2-10} alkenyl group, \underline{or} a C_{2-10} alkenyl group \underline{or} an \underline{aryl} group), a group represented by formula \underline{fi}

$$\mathbb{R}^{d}$$
 \mathbb{O}

(wherein R⁴ is the same as described above) or a group represented by formula [ii]

36. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a C₁₋₁₀alkyl group, a compositive d by one or two aryl groups, a hydroxyC₂₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkyl group or a C₁₋₁₀alkyl group, a farnesyl group, a 4-morpholinylC₁₋₁₀alkyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group).

 (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim

2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -SCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

R¹ represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond, R^c represents a hydrogen atom, C₁.

10alkyl group, or a C₂₋₁₀alkenyl group-or-an aryl-group; and R^d represents a C₁₋₁₀alkyl group; or a

C₂₋₁₀alkenyl group-or-an aryl-group), a group represented by formula fil

$$\mathbb{R}^{d}$$
 \mathbb{O}

(wherein R^d is the same as described above) or a group represented by formula [ii]

38. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -SR³ (wherein R³ is the same as described above); and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a C₁₋₁₀alkyl group substituted by one or two aryl groups, a hydroxyC₂₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkyl group, a

₁₀alkoxycarbonylC₁₋₁₀alkyl group, a farnesyl group, a 4 morpholinylC₁₋₁₀alkyl group, or a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR*R^b (wherein R* and R* are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group).

39. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents-SR³ (wherein R³ is the same as described above); and

 R^1 represents a group represented by formula-CHR c OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or an aryl group; and R^d represents a C_{1-10} alkyl group; or a C_{2-10} alkenyl group-or an aryl group, a group represented by formula [i]

(wherein R^d is the same as described above) or a group represented by formula [ii]

 (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim

2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -S(O), CHR³R⁴ (wherein R³, R⁴ and n are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

41. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -S(O)_nCHR³R⁴ (wherein R³, R⁴ and n are the same as described above); and

 R^1 represents a group represented by formula-CHR OC(O)ZR (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, $C_{1.0}$ alkyl group, or a $C_{2.10}$ alkenyl group or an aryl group; and R° represents a $C_{1.10}$ alkyl group, or a $C_{2.10}$ alkenyl group, a group represented by formula [i]

$$\mathbb{R}^{d}$$
 \mathbb{O}

(wherein R^d is the same as described above) or a group represented by formula [ii]

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42. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{2-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

43. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate-thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -NHCHR³R⁴ (wherein R³ and R⁴ are the same as described above); and

R¹ represents a group represented by formula-CHR°OC(O)XR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, a C₁₋₁₀alkyl group, or a C₂₋₁₀alkenyl group or an aryl group; and R^d represents a C₁₋₁₀alkyl group; or a C₂₋₁₀alkenyl group, a group represented by formula fil

$$\mathbb{R}^{d}$$
 \mathbb{O}

(wherein R^d is the same as described above) or a group represented by formula [ii]

44. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -N(CHR³R⁴)(CHR³'R⁴) (wherein R³, R³, R⁴ and R⁴ are the same as described above); and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a-C₁₋₁₀alkyl group, a balogenoC₁₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, a nazidoC₁₋₁₀alkyl group, a naminoC₂₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxyCarbonylC₁₋₁₀alkyl group, a farnesyl group, a 4-morpholinylC₁₋₁₀alkyl group or a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group).

45. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative. or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim

2, wherein in the formula [II], R^2 represents a hydrogen atom; X represents a hydrogen atom; Y represents -N(CHR³R⁴)(CHR³'R^{4'}) (wherein R^3 , R^3 ', R^4 and $R^{4'}$ are the same as described above); and

 R^1 represents a group represented by formula-CHR $^{\circ}$ OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, $C_{1.00}$ alkyl group, \underline{or} a $C_{2.10}$ alkenyl group-or an aryl group; and R^d represents a $C_{1.10}$ alkyl group; \underline{or} a $C_{2.10}$ alkenyl group, a group represented by formula [i]

(wherein R^d is the same as described above) or a group represented by formula [ii]

46. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ represents a hydrogen atom; R⁴ represents a phenyl group or a phenyl group substituted by one to five substituents selected from a group coensisting of a halogen atom, a phenyl group, a C_{1.10}alkyl group, a C_{1.10}alkoxy group, a trifluoromethyl group, a

phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, a nazido C_{1-10} alkyl group, a namino C_{2-10} alkyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

47. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ represents a hydrogen atom; R⁴ represents a phenyl group or a phenyl group substituted by one to five substitutents substituents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a group represented by formula-CHR $^{\circ}$ OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, a C_{1-10} alkyl group, or a C_{2-10} alkenyl group-or-an-aryl group; and R^d represents a C_{1-10} alkyl group, or a C_{2-10} alkenyl group or an aryl group, a group represented by formula [i]

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$$\mathbb{R}^{d}$$
 O O

(wherein R^d is the same as described above) or a group represented by formula [ii]

- 48. (canceled).
- 49. (canceled).
- 50. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are identical or different, and each represents a phenyl group or a phenyl group substituted by one to five substitutents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkyl group, a C

 $_{10}$ alkoxycarbonyl C_{1-10} alkyl group, a farnesyl group, a 4-morpholinyl C_{1-10} alkyl group or a C_1 . $_{10}$ alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

51. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a fluorine atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are identical or different, and each represents a phenyl group or a phenyl group substituted by one to five substitutents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a group represented by formula-CHR $^{\circ}$ OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R° represents a hydrogen atom, a C_{1-10} alkyl group, \underline{or} a C_{2-10} alkenyl group- \underline{or} an \underline{aryl} group; and R^d represents a C_{1-10} alkyl group, \underline{or} a C_{2-10} alkenyl group \underline{or} and \underline{or} aryl group, \underline{or} a \underline{or} a \underline{or} group represented by formula [i]

$$\mathbb{R}^{d}$$
 \mathbb{Q}
 \mathbb{Q}

(wherein R^d is the same as described above) or a group represented by formula [ii]

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52. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [III], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -OCHR³R⁴ (wherein R³ represents a hydrogen atom; R⁴ represents a phenyl group or a phenyl group substituted by one to five substitutents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

R¹ represents a C₁₋₁₀alkyl group, a C₂₋₁₀alkenyl group, a C₂₋₁₀alkynyl group, a -C₁₋₁₀alkyl group, a -C₁₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, a halogenoC₁₋₁₀alkyl group, an azidoC₁₋₁₀alkyl group, an aminoC₂₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxyC₁₋₁₀alkyl group, a C₁₋₁₀alkoxyCarbonylC₁₋₁₀alkyl group, a farnesyl group, a 4-morpholinylC₁₋₁₀alkyl group or a C₁₋₁₀alkyl group substituted by a group represented by formula-C(O)NR^aR^b (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C₁₋₁₀alkyl group).

53. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents a COCHR³R⁴ (wherein R³ represents a hydrogen atom; R⁴ represents a phenyl group or a

phenyl group substituted by one to five substituents substituents selected from a group containing a halogen atom, a phenyl group, a C_{1-10} alkyl group, a C_{1-10} alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and phenoxy group); and

 R^1 represents a group represented by formula-CHR c OC(O)ZR d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C_{1-10} alkyl group, \underline{or} a C_{2-10} alkenyl group- \underline{or} and R^d represents a C_{1-10} alkyl group- \underline{or} a C_{2-10} alkenyl group- \underline{or} a C_{2-10} alkenyl group- \underline{or} a C_{2-10} alkenyl group- \underline{or} and \underline{or} a C_{2-10} alkenyl group- \underline{or} and \underline{or} a \underline{or} and \underline{or} a \underline{or} and \underline{or} a \underline{or} a \underline{or} and \underline{or} a \underline{or} and \underline{or} an

(wherein R4 is the same as described above) or a group represented by formula [ii]

- 54. (canceled).
- 55. (canceled).
- **56.** (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaceutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y

represents -OCHR³R⁴ (wherein R³ and R⁴ are identical or different, and each represents a phenyl group or a phenyl group substituted by one to five substitutents-substituents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

 R^1 represents a C_{1-10} alkyl group, a C_{2-10} alkenyl group, a C_{2-10} alkynyl group, a C_{1-10} alkyl group, a halogeno C_{1-10} alkyl group, an azido C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group, an amino C_{2-10} alkyl group, a C_{1-10} alkyl group, a C_{1-10} alkyl group, a farnesyl group, a C_{1-10} alkyl group or a C_{1-10} alkyl group substituted by a group represented by formula- $C(O)NR^aR^b$ (wherein R^a and R^b are identical or different, and each represents a hydrogen atom or a C_{1-10} alkyl group).

57. (currently amended): A 2-amino-bicyclo[3.1.0]hexane-2,6-dicarboxylic ester derivative, or a pharmaccutically acceptable salt thereof or a hydrate thereof-according to claim 2, wherein in the formula [II], R² represents a hydrogen atom; X represents a hydrogen atom; Y represents -OCHR³R⁴ (wherein R³ and R⁴ are identical or different, and each represents a phenyl group or a phenyl group substituted by one to five substituents-substituents selected from a group containing a halogen atom, a phenyl group, a C₁₋₁₀alkyl group, a C₁₋₁₀alkoxy group, a trifluoromethyl group, a phenyl group, a hydroxycarbonyl group, an amino group, a nitro group, a cyano group and a phenoxy group); and

R¹ represents a group represented by formula-CHR^cOC(O)ZR^d (wherein Z represents an oxygen atom, a nitrogen atom, a sulfur atom or a single bond; R^c represents a hydrogen atom, a C₁₋₁₀alkyl group, <u>or a C₂₋₁₀alkenyl group or an aryl group;</u> and R^d represents a C₁₋₁₀alkyl group; or a C₂₋₁₀alkenyl group or an aryl group represented by formula fil

$$\mathbb{R}^{d}$$
 0 0 0

(wherein R^d is the same as described above) or a group represented by formula [ii]

- **58.** (currently amended): A drug comprising the 2-amino-bicyclo [3.1.0] hexane 2,6-dicarboxylic ester-derivative, or the pharmaceutically acceptable salt thereof or the hydrate thereof according claim 2 as an active ingredient.
- 59. (original): A drug according to claim 58, wherein the drag is a group II metabotropic glutamate receptor antagonist.
- 60. (previously presented): (1R,2R,3R,5R,6R)-2-amino-3-(3,4-dichlorobenzyloxy)-6-fluoro-2,6-dicarboxylic acid 6-n-heptyl ester represented by the following structure: